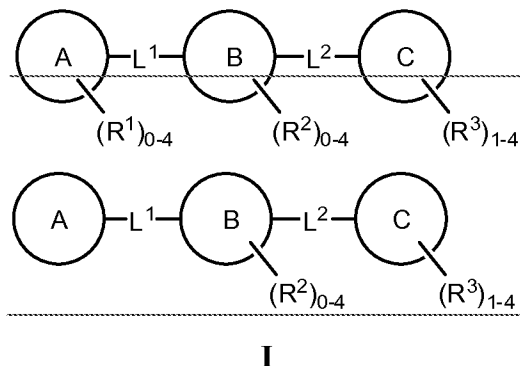


Amendments to the Claims

The following listing of claims will replace all prior versions and listings of claims in the application.

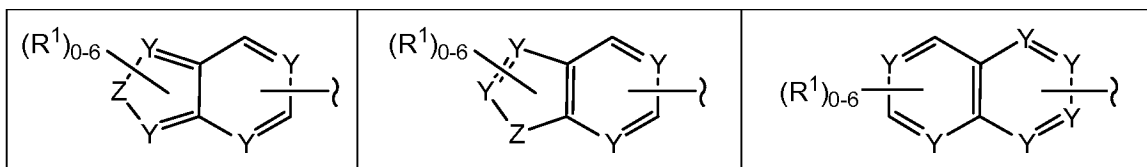
Listing of Claims:

1. (Currently amended) A compound for modulating c-Kit activity according to Formula I,



or a pharmaceutically acceptable salt, thereof, wherein,

ring A is:



wherein each Y is independently either =C(H)- or =N-; and Z is selected from -O-, -S-, and -N(R⁷)-, provided that the A ring contains at least one annular N, O, or S;

each R¹ is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two adjacent of R¹, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R¹⁰;

L¹ is a single bond;

ring B is phenyl;

each R² is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴, -N(R⁴)R⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two adjacent of R², together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R¹⁵;

L² is selected from -N(H)N(H)C(=O)N(H)-, -CH₂N(H)C(=O)N(H)-, -CH₂OC(=O)N(H)-, and -XCH₂C(=O)N(H)-; wherein X is selected from -O-, -S(O)₀₋₂-, and -N(R⁷)-; and any C-H of L² is optionally C-R²⁰;

ring C is phenyl;

each R^3 is independently selected from halogen, trihalomethyl, $-\text{CN}$, $-\text{NO}_2$, $-\text{OR}^4$, $-\text{N}(\text{R}^4)\text{R}^4$, $-\text{S}(\text{O})_{0-2}\text{R}^4$, $-\text{SO}_2\text{N}(\text{R}^4)\text{R}^4$, $-\text{C}(=\text{O})\text{N}(\text{R}^4)\text{R}^4$, $-\text{C}(=\text{NR}^5)\text{N}(\text{R}^4)\text{R}^4$, $-\text{C}(=\text{NR}^5)\text{R}^4$, $-\text{N}(\text{R}^4)\text{SO}_2\text{R}^4$, $-\text{N}(\text{R}^4)\text{C}(\text{O})\text{R}^4$, $-\text{NCO}_2\text{R}^4$, $-\text{C}(=\text{O})\text{R}^4$, optionally substituted alkoxy, optionally substituted C_{1-6} alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6} alkyl, ~~optionally substituted heterocyclyl,~~ and optionally substituted heterocyclyl C_{1-6} alkyl; provided R^3 is not a cyclic sulfonamide attached to ring C via the nitrogen of said cyclic sulfonamide,

wherein there exists at least one of R^3 that is halogen or trihalomethyl;

R^4 is selected from $-\text{H}$, optionally substituted C_{1-6} alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6} alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6} alkyl;

two of R^4 , when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R^5 is selected from $-\text{H}$, $-\text{CN}$, $-\text{NO}_2$, $-\text{OR}^4$, $-\text{S}(\text{O})_{0-2}\text{R}^4$, $-\text{CO}_2\text{R}^4$, optionally substituted C_{1-6} alkyl, optionally substituted C_{1-6} alkenyl, and optionally substituted C_{1-6} alkynyl;

R^7 is selected from $-\text{H}$, optionally substituted C_{1-6} alkyl, $-\text{SO}_2\text{N}(\text{R}^4)\text{R}^4$, $-\text{CO}_2\text{R}^4$, $-\text{C}(=\text{O})\text{N}(\text{R}^4)\text{R}^4$, $-\text{C}(=\text{NR}^5)\text{N}(\text{R}^4)\text{R}^4$, $-\text{C}(=\text{NR}^5)\text{R}^4$, $-\text{C}(=\text{O})\text{R}^4$, optionally substituted alkoxy, optionally substituted aryl C_{1-6} alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6} alkyl; and

each of R^{10} , each of R^{15} , each of R^{20} , and each of R^{25} is independently selected from $-\text{H}$, halogen, trihalomethyl, $-\text{CN}$, $-\text{NO}_2$, $-\text{OR}^4$, $-\text{N}(\text{R}^4)\text{R}^4$, $-\text{S}(\text{O})_{0-2}\text{R}^4$, $-\text{SO}_2\text{N}(\text{R}^4)\text{R}^4$, $-\text{CO}_2\text{R}^4$, $-\text{C}(=\text{O})\text{N}(\text{R}^4)\text{R}^4$, $-\text{C}(=\text{NR}^5)\text{N}(\text{R}^4)\text{R}^4$, $-\text{C}(=\text{NR}^5)\text{R}^4$, $-\text{N}(\text{R}^4)\text{SO}_2\text{R}^4$, $-\text{N}(\text{R}^4)\text{C}(\text{O})\text{R}^4$, $-\text{NCO}_2\text{R}^4$, $-\text{C}(=\text{O})\text{R}^4$, optionally substituted alkoxy, optionally substituted C_{1-6} alkyl, optionally substituted aryl, optionally substituted aryl

C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

provided:

the compound is not one of:

N-(2,3-dichlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,

N-(4-chloro-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,

N-(4-bromophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide

N-(2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide

N-(4-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide

2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoromethyl)phenyl]acetamide

2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl)phenyl]acetamide

N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide

N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy}acetamide or

N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide.

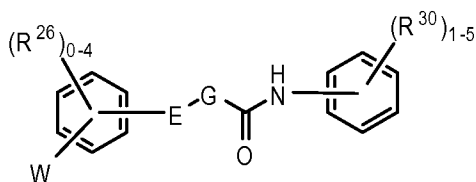
2-12. (Cancelled)

13. (Canceled)

14. (Original) The compound according to claim 13, wherein ring C is a phenyl comprising a trifluoromethyl radical *meta*- to L².

15. (Previously presented) The compound according to claim 1, wherein each of R³ is independently selected from halogen, trihalomethyl, -OR⁴, -C(=O)R⁴, and optionally substituted C₁₋₆alkyl.

16. (Currently amended) A compound for modulating c-Kit activity according to the following Formula:



II

or a pharmaceutically acceptable salt, thereof, wherein,

W is selected from the following:

each of R^{27} independently selected from halogen, trihalomethyl, $-\text{CN}$, $-\text{NO}_2$, $-\text{OR}^{55}$, $-\text{S}(\text{O})_{0-2}\text{R}^{55}$, $-\text{SO}_2\text{N}(\text{R}^{55})\text{R}^{55}$, $-\text{C}(=\text{O})\text{N}(\text{R}^{55})\text{R}^{55}$, $-\text{C}(=\text{NR}^{50})\text{N}(\text{R}^{55})\text{R}^{55}$, $-\text{C}(=\text{NR}^{50})\text{R}^{55}$, $-\text{N}(\text{R}^{55})\text{SO}_2\text{R}^{55}$, $-\text{N}(\text{R}^{55})\text{C}(\text{O})\text{R}^{55}$, $-\text{NCO}_2\text{R}^{55}$, $-\text{C}(=\text{O})\text{R}^{55}$, optionally substituted alkoxy, optionally substituted C_{1-6} alkyl, optionally substituted aryl C_{1-6} alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6} alkyl;

each Y is independently either $=\text{C}(\text{H})-$ or $=\text{N}-$;

Z is selected from -O-, -S(O)₀₋₂-, and -N(R⁷)-,

provided that the W ring contains at least one annular N, O, or S;

E and G are each independently selected from -O-, -S(O)₀₋₂-, -C(R³¹)R³²-, and -N(R³³)-,

J₁ and J₂ are each independently =C(H)- or =N-;

R²⁶ is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -S(O)₀₋₂R⁴⁰, -SO₂N(R⁴⁰)R⁴⁰, -CO₂R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -N(R⁴⁰)SO₂R⁴⁰, -N(R⁴⁰)C(O)R⁴⁰, -NCO₂R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

R³⁰ is independently selected from halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -S(O)₀₋₂R⁴⁰, -SO₂N(R⁴⁰)R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -N(R⁴⁰)SO₂R⁴⁰, -N(R⁴⁰)C(O)R⁴⁰, -NCO₂R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, ~~optionally substituted heterocyclyl,~~ and optionally substituted heterocyclyl C₁₋₆alkyl, wherein there exists at least one of R³⁰ that is trihalomethyl;

R³¹ and R³² are each independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -S(O)₀₋₂R⁴⁰, -SO₂N(R⁴⁰)R⁴⁰, -CO₂R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -N(R⁴⁰)SO₂R⁴⁰, -N(R⁴⁰)C(O)R⁴⁰, -NCO₂R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

R³³ is selected from -H, optionally substituted lower alkyl, -SO₂N(R⁴⁰)R⁴⁰, -CO₂R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

R⁴⁰ is selected from -H, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two of R⁴⁰, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R⁵⁰ is selected from -H, -CN, -NO₂, -OR⁴⁰, -S(O)₀₋₂R⁴⁰, -CO₂R⁴⁰, optionally substituted C₁₋₆alkyl, optionally substituted C₁₋₆alkenyl, and optionally substituted C₁₋₆alkynyl;

R⁵⁵ is selected from -H, optionally substituted C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl; and

two of R⁵⁵, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P.

17. (Cancelled)

18. (Currently amended) The compound according to claim 16, wherein R³⁰ is selected from halogen, trihalomethyl, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, ~~optionally substituted heterocyclyl,~~ and optionally substituted heterocyclyl C₁₋₆alkyl, wherein there exists at least one of R³⁰ that is trifluoromethyl.

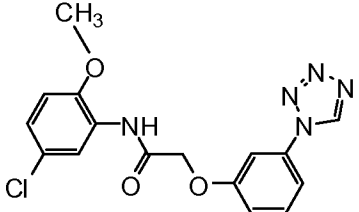
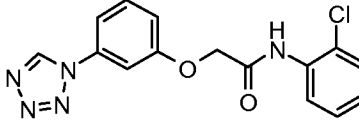
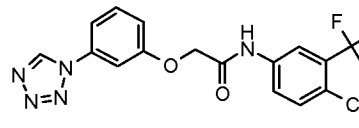
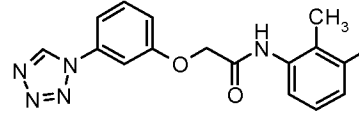
19. (Cancelled)

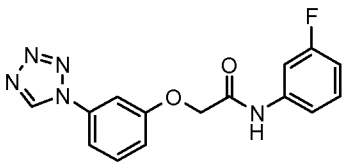
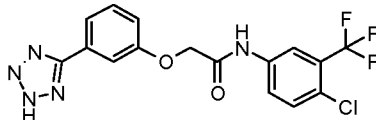
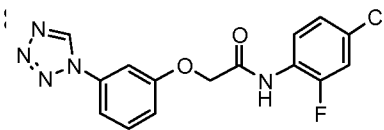
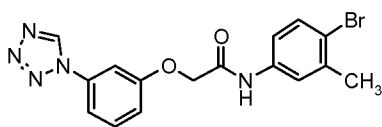
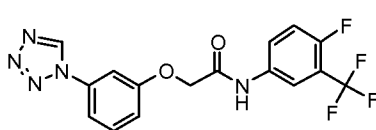
20. (Cancelled)

21. (Cancelled)

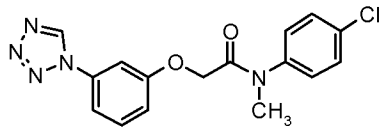
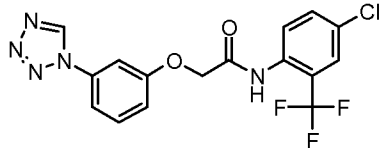
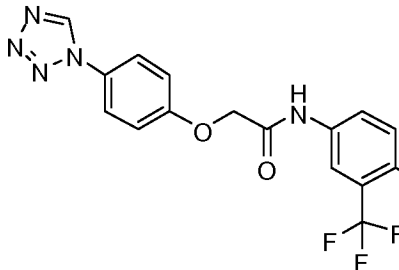
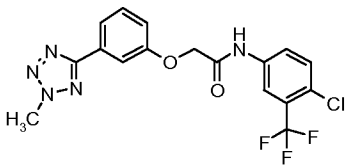
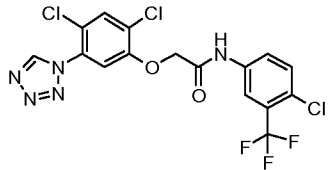
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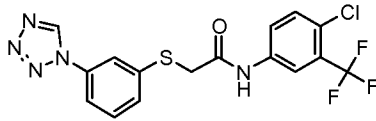
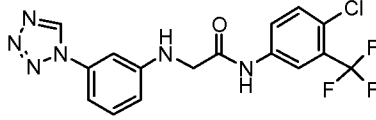
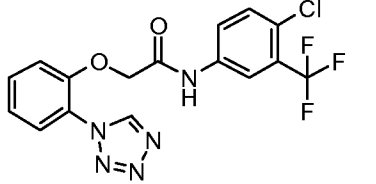
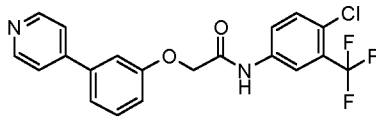
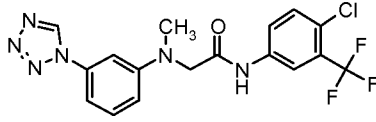
23. (Withdrawn) The compound according to claim 16, wherein E is selected from -O-, -S(O)₀₋₂-, and -NH-; and G is -CH₂-.
24. (Withdrawn) The compound according to claim 16, wherein E is either -CH₂- or -NH-; and G is selected from -O-, -S-, and -NH-.
25. (Cancelled)
26. (Cancelled)
27. (Previously presented) A compound selected from the following Table:

Entry	Name	Structure
1	N-[5-chloro-2-(methoxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
4	N-(2-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
5	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
7	N-(3-chloro-2-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Entry	Name	Structure
8	N-(3-fluorophenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
9	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(2H-tetrazol-5-yl)phenyl]oxy} acetamide	
10	N-(4-chloro-2-fluorophenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
11	N-(4-bromo-3-methylphenyl)-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
13	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	

Entry	Name	Structure
14	N-[4-bromo-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]oxy} acetamide	
15	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[4-(1H-tetrazol-1- yl)phenyl]oxy} acetamide	
16	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]oxy} propanamide	
17	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(5-methyl-1H-tetrazol-1- yl)phenyl]oxy} acetamide	
18	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2-methyl-5-(1H-tetrazol-1- yl)phenyl]oxy} acetamide	

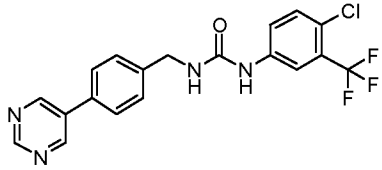
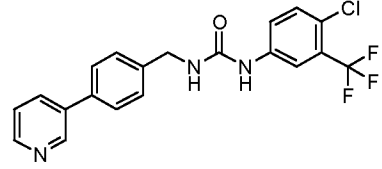
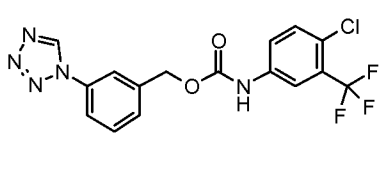
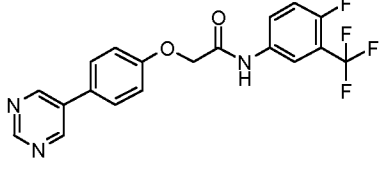
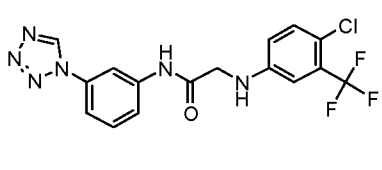
Entry	Name	Structure
19	N-(4-chlorophenyl)-N-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
20	N-[4-chloro-2-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
23	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	
24	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(2-methyl-2H-tetrazol-5-yl)phenyl]oxy}acetamide	
25	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2,4-dichloro-5-(1H-tetrazol-1-yl)phenyl]oxy}acetamide	

Entry	Name	Structure
26	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]thio} acetamide	
27	N-[4-chloro-3-(trifluoromethyl)phenyl]- N~2~-[3-(1H-tetrazol-1- yl)phenyl]glycinamide	
28	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2-(1H-tetrazol-1- yl)phenyl]oxy} acetamide	
37	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-pyridin-4-ylphenyl)oxy]acetamide	
38	N-[4-chloro-3-(trifluoromethyl)phenyl]- N~2~-methyl-N~2~-[3-(1H-tetrazol-1- yl)phenyl]glycinamide	

Entry	Name	Structure
49	N-[5-chloro-2,4-bis(methoxy)phenyl]-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
52	N-[2-(methoxy)-5-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
57	1,1-dimethylethyl 2-{3-[(2-{[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl)oxy]phenyl}-1H-pyrrole-1-carboxylate	
58	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-pyrrol-2-yl)phenyl]oxy} acetamide	
59	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-pyrimidin-5-ylphenyl)oxy]acetamide	

Entry	Name	Structure
60	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-1,2,3-triazol-1- yl)phenyl]oxy} acetamide	
63	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-pyridin-3-ylphenyl)oxy]acetamide	
68	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(3,5-dimethylisoxazol-4- yl)phenyl]oxy} acetamide	
69	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-quinolin-7-ylphenyl)oxy]acetamide	
71	N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [3-(1H-tetrazol-1- yl)phenyl]hydrazinecarboxamide	

Entry	Name	Structure
73	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	
75	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-tetrazol-1-yl)phenyl]methyl}urea	
77	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N~2~-[3-(1H-tetrazol-1-yl)phenyl]glycinamide	
79	N-[2-fluoro-5-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
80	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide	
81	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyrimidin-5-ylphenyl)methyl]urea	

Entry	Name	Structure
82	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyrimidin-5-ylphenyl)methyl]urea	
83	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	
84	[3-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
85	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide	
86	N-[3-(1H-tetrazol-1-yl)phenyl]-N'-[4-chloro-3-(trifluoromethyl)phenyl]glycinamide	

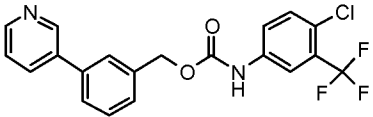
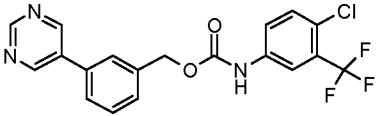
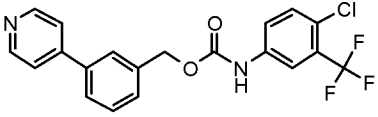
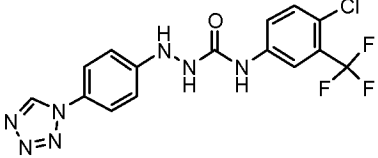
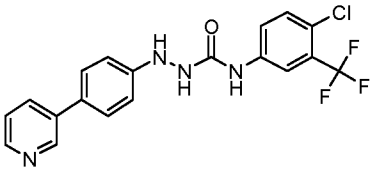
Entry	Name	Structure
87	2-{[4-chloro-3-(trifluoromethyl)phenyl]oxy}-N-[3-(1H-tetrazol-1-yl)phenyl]acetamide	
88	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-methyl-4-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
89	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-1,2,3-triazol-1-yl)phenyl]oxy} acetamide	
90	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-fluoro-4-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	
91	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2-fluoro-4-(1H-tetrazol-1-yl)phenyl]oxy} acetamide	

Entry	Name	Structure
94	N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide	
95	2-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-fluoro-3-(trifluoromethyl)phenyl]acetamide	
96	2-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}oxy)-N-[4-chloro-3-(trifluoromethyl)phenyl]acetamide	
97	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-4-ylphenyl)oxy]acetamide	
98	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(methyloxy)-4-(1H-tetrazol-1-yl)phenyl]glycinamide	

Entry	Name	Structure
99	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[4-(methoxy)-3-(1H-tetrazol-1-yl)phenyl]glycinamide	
100	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[4-(1H-tetrazol-1-yl)phenyl]glycinamide	
101	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(2,3,5,6-tetrafluoro-4-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
102	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(1H-tetrazol-1-yl)phenyl]methyl}urea	
103	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyrimidin-5-ylphenyl)hydrazinecarboxamide	

Entry	Name	Structure
104	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyridin-3-ylphenyl)methyl]urea	
105	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
106	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}propanamide	
107	N-({4-[2,4-bis(methoxy)pyrimidin-5-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
108	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[2-(methoxy)pyrimidin-5-yl]phenyl}methyl)urea	

Entry	Name	Structure
109	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[6-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
110	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[2-(methyloxy)pyrimidin-5-yl]phenyl}methyl)urea	
111	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[6-(methyloxy)pyridin-3-yl]phenyl}methyl)urea	
114	N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-([3-(2H-tetrazol-5-yl)phenyl]glycinamide	
115	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{2,6-difluoro-4-(1H-tetrazol-1-yl)phenyl}oxy}acetamide	

Entry	Name	Structure
116	(3-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
117	(3-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
118	(3-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
119	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[4-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide	
120	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-3-ylphenyl)hydrazinecarboxamide	

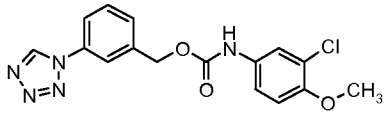
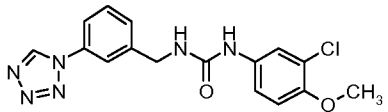
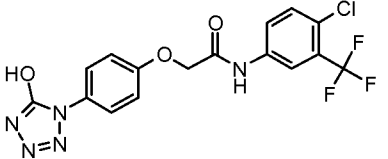
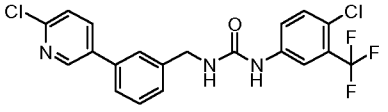
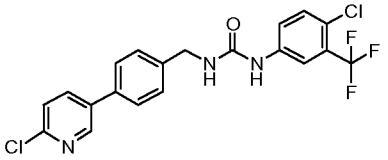
Entry	Name	Structure
121	(4-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
122	(4-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
123	(4-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
124	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-4-ylphenyl)methyl]urea	
125	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-3-ylphenyl)hydrazinecarboxamide	

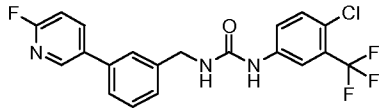
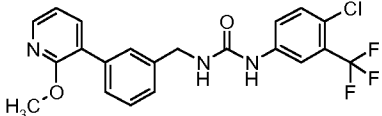
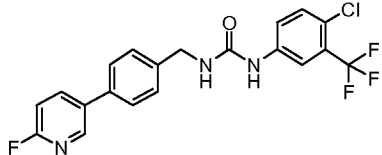
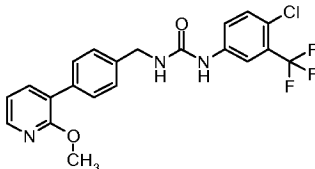
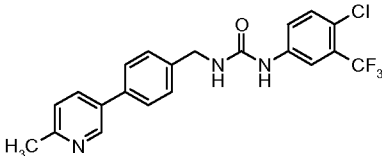
Entry	Name	Structure
126	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide	
127	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(4-pyrimidin-5-ylphenyl)methyl]urea	
128	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	
129	(4-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	
130	(4-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	

Entry	Name	Structure
131	1-(4-pyridin-3-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
132	1-(4-pyrimidin-5-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
133	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(3-pyridin-3-ylphenyl)methyl]urea	
134	N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(3-pyrimidin-5-ylphenyl)methyl]urea	
135	(3-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	

Entry	Name	Structure
136	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate	
138	N-[4-fluoro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea	
143	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyridin-3-ylphenyl)ethyl]urea	
144	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyrimidin-5-ylphenyl)ethyl]urea	
145	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[[4-(1H-indol-2-yl)phenyl]oxy]acetamide	

Entry	Name	Structure
147	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-4-ylphenyl)hydrazinecarboxamide	
148	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-4-ylphenyl)hydrazinecarboxamide	
149	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyridin-4-ylphenyl)methyl]urea	
150	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-quinoxalin-6-ylphenyl)methyl]urea	
152	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-quinoxalin-6-ylphenyl)methyl]urea	

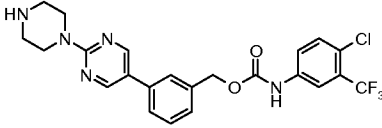
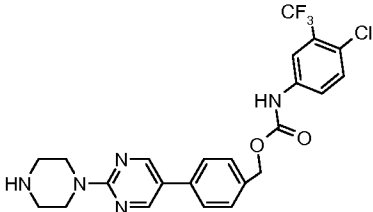
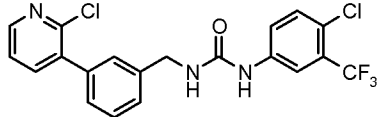
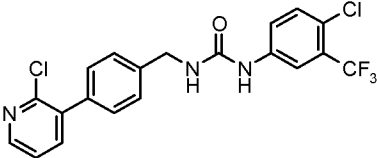
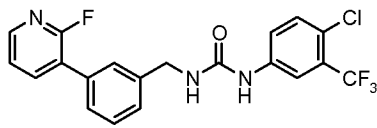
Entry	Name	Structure
155	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-chloro-4-(methyloxy)phenyl]carbamate	
156	N-[3-chloro-4-(methyloxy)phenyl]-N'-{[3-(1H-tetrazol-1-yl)phenyl]methyl}urea	
157	N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(5-hydroxy-1H-tetrazol-1-yl)phenyl]oxy}acetamide	
160	N-{[3-(6-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
161	N-{[4-(6-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	

Entry	Name	Structure
165	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[3-(6-fluoropyridin-3-yl)phenyl]methyl}urea	
166	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({3-[2-(methoxy)pyridin-3-yl]phenyl}methyl)urea	
167	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(6-fluoropyridin-3-yl)phenyl]methyl}urea	
168	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({4-[2-(methoxy)pyridin-3-yl]phenyl}methyl)urea	
169	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(6-methylpyridin-3-yl)phenyl]methyl}urea	

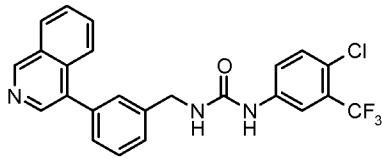
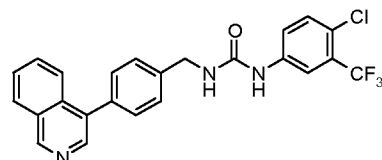
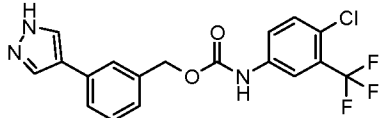
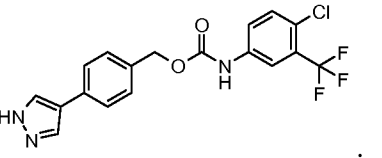
Entry	Name	Structure
171	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(6-methylpyridin-3-yl)phenyl]methyl}urea	
174	[3-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
177	(3-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
179	N-{[3-(6-acetylpyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
180	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(6-cyanopyridin-3-yl)phenyl]methyl}urea	
185	[3-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Entry	Name	Structure
188	[3-(1H-benzimidazol-2-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
190	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({3-[5-(methylthio)pyridin-3-yl]phenyl}methyl)urea	
191	[4-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
194	(4-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
195	[4-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Entry	Name	Structure
200	(3-pyridin-3-ylphenyl)methyl [5-chloro-2-(methyloxy)phenyl]carbamate	
201	[4-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
202	(3-pyrimidin-5-ylphenyl)methyl [5-chloro-2-(methyloxy)phenyl]carbamate	
213	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-pyrazol-4-yl)phenyl]methyl} urea	
214	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(1H-pyrazol-4-yl)phenyl]methyl} urea	

Entry	Name	Structure
215	[3-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
216	[4-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
217	N-{[3-(2-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
218	N-{[4-(2-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea	
219	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(2-fluoropyridin-3-yl)phenyl]methyl}urea	

Entry	Name	Structure
220	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(2-fluoropyridin-3-yl)phenyl]methyl}urea	
221	[3-(1H-tetrazol-1-yl)phenyl]methyl [3-(trifluoromethyl)phenyl]carbamate	
224	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[5-(methylthio)pyridin-2-yl]phenyl}methyl)urea	
225	[3-(2,6-dimethylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
226	{3-[5-(methyloxy)pyridin-3-yl]phenyl}methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

Entry	Name	Structure
229	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-isoquinolin-4-ylphenyl)methyl]urea	
230	N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-isoquinolin-4-ylphenyl)methyl]urea	
232	[3-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	
233	[4-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate	

28. (Previously presented) A pharmaceutical composition comprising the compound according to claim 1 and a pharmaceutically acceptable carrier.
29. (Cancelled)
30. (Withdrawn) A method for modulating the *in-vivo* activity of a kinase, the method comprising administering to a subject an effective amount of the compound according to claim 1.
31. (Withdrawn) The method according to claim 30, wherein the kinase is c-Kit.

32. (Withdrawn) The method according to claim 31, wherein modulating the *in vivo* activity of c-Kit comprises inhibition of c-Kit.
33. (Withdrawn) A method of treating rheumatoid arthritis, graft-host diseases, multiple sclerosis, psoriasis; arterosclerosis, myocardioinfarction, ischemia, stroke, restenosis; interbowel diseases, osteoarthritis, macular degeneration, or diabetic retinopathy, the method comprising administering, to a mammal in need thereof, a therapeutically effective amount of the compound or the pharmaceutical composition as described in claim 1 .
34. (Withdrawn) A method of screening for modulators of c-Kit, the method comprising combining the compound according to claim 1 and at least one candidate agent and determining the effect of the candidate agent on c-Kit activity.
35. (Withdrawn) A method of inhibiting proliferative activity in a cell, the method comprising administering an effective amount of a composition comprising the compound according to claim 1 to a cell or a plurality of cells.